

NEWS 7 Apr 22 BIOSIS Gene Names now available in TOXCENTER
 NEWS 8 Apr 22 Federal Research in Progress (FEDRIP) now available
 NEWS 9 Jun 03 New e-mail delivery for search results now available
 NEWS 10 Jun 10 MEDLINE Reload
 NEWS 11 Jun 10 PCTFULL has been reloaded
 NEWS 12 Jul 02 FOREGE no longer contains STANDARDS file segment
 NEWS 13 Jul 22 USAN to be reloaded July 28, 2002;
 saved answer sets no longer valid
 NEWS 14 Jul 29 Enhanced polymer searching in REGISTRY
 NEWS 15 Jul 30 NETFIRST to be removed from STN
 NEWS 16 Aug 08 CANCERLIT reload
 NEWS 17 Aug 08 PHARMAMarketLetter(PHARMAML) - new on STN
 NEWS 18 Aug 08 NTIS has been reloaded and enhanced
 NEWS 19 Aug 09 JAPIO to be reloaded August 25, 2002
 NEWS 20 Aug 19 Aquatic Toxicity Information Retrieval (AQUIRE)
 now available on STN
 NEWS 21 Aug 19 IFIPAT, IFICDB, and IFIUDB have been reloaded
 NEWS 22 Aug 19 The MEDLINE file segment of TOXCENTER has been reloaded
 NEWS 23 Aug 26 Sequence searching in REGISTRY enhanced

 NEWS EXPRESS February 1 CURRENT WINDOWS VERSION IS V6.0d,
 CURRENT MACINTOSH VERSION IS V6.0a(ENG) AND V6.0Ja(JP),
 AND CURRENT DISCOVER FILE IS DATED 05 FEBRUARY 2002
 NEWS HOURS STN Operating Hours Plus Help Desk Availability
 NEWS INTER General Internet Information
 NEWS LOGIN Welcome Banner and News Items
 NEWS PHONE Direct Dial and Telecommunication Network Access to STN
 NEWS WWW CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that specific topic.

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* * * * * STN Columbus * * * * *

FILE 'HOME' ENTERED AT 06:22:05 ON 28 AUG 2002

=> file cacplus

'CACPLUS' IS NOT A VALID FILE NAME

SESSION CONTINUES IN FILE 'HOME'

Enter "HELP FILE NAMES" at an arrow prompt (=>) for a list of files that are available. If you have requested multiple files, you can specify a corrected file name or you can enter "IGNORE" to continue accessing the remaining file names entered.

=> e

N-[N-[3-(3-methoxy-4-hydroxyphenyl)propyl]-L-.alpha.-aspartyl]-L-phenylalanine
1-Me ester/cn

THIS COMMAND NOT AVAILABLE IN THE CURRENT FILE

The EXPAND command is used to look at the index in a file which has an index. This file does not have an index.

=> file reg
COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
0.63	0.63

FULL ESTIMATED COST

FILE 'REGISTRY' ENTERED AT 06:23:52 ON 28 AUG 2002
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STRUCTURE FILE UPDATES: 26 AUG 2002 HIGHEST RN 444986-65-6
DICTIONARY FILE UPDATES: 26 AUG 2002 HIGHEST RN 444986-65-6

TSCA INFORMATION NOW CURRENT THROUGH MAY 20, 2002

Please note that search-term pricing does apply when
conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Calculated physical property data is now available. See HELP PROPERTIES
for more information. See STNote 27, Searching Properties in the CAS
Registry File, for complete details:
<http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf>

=> e

N-[N-[3-(3-methoxy-4-hydroxyphenyl)propyl]-L-.alpha.-aspartyl]-L-phenylalanine
1-Me ester/cn

E1	1	N,S-DIVINYL-2-MERCAPTOBENZIMIDAZOLE/CN
E2	1	N,S-TRANS-1,2-DITHIOLANYLIDENETHIAMINE/CN
E3	0 -->	N-N-3-(3-METHOXY-4-HYDROXYPHENYL) PROPYL -L-.ALPHA.-ASPARTY L -L-PHENYLALANINE 1-ME ESTER/CN

E4 1
N-(((4-(2-(5-CHLORO-2-METHOXYBENZOYL)(4-NITROPHENYL) SULFO

NYL)AMINO)ETHYL)PHENYL)SULFONYL)(4-NITROPHENYL)SULFONYL)AMI
NO)CARBONYL)-N-CYCLOHEXYL-4-NITROBENZENESULFONAMIDE/CN

E5 1
N-(((1-(3,4-DIHYDROXY-5-(HYDROXYMETHYL)TETRAHYDRO-2-FURANYL)

-5-iodo-2-oxo-1,2-dihydro-4-pyrimidinyl)(4-nitrophenyl) sulf

ONYL)AMINO)CARBOTHIOYL)-4-NITRO-N-(4-NITROPHENYL)BENZENESULF
ONAMIDE/CN

E6	1	N-((2,3-DIMETHOXYBENZYL)OXY)CARBONYL)TRYPTOPHAN/CN
E7	1	N-((2,6-DICHLOROPHENYL)CARBAMOYL)METHYL)IMINODIACETIC ACID/

E8	1	CN N-((2,6-DIISOPROPYLPHENYL)CARBAMOYL)METHYL)IMINODIACETIC AC
----	---	--

E9	1	ID/CN N-((2-(DIETHYLAMINO)ETHYL)OXY)CARBONYL)CYCLOPEPTIDE A, HYDR
----	---	---

E10	1	OCHLORIDE/CN N-((2-AMINOETHYL)AMINOMETHYL)BENZYL)DIMETHYLSILANE/CN
E11	1	N-((2-HYDROXYBENZOYL)OXY)METHYL)BENZAMIDE/CN
E12	1	N-((3,4-DIMETHOXYPHENYL)METHYLENE)AMINO)PHTHALIMIDE/CN

=> e

N-(N-(3-(3-methoxy-4-hydroxyphenyl)propyl)-L-.alpha.-aspartyl)-L-phenylalanine
1-Me ester/cn

E1 1 N-(N-(2-METHYLPHENYL)-N-PHENYLAMINO) PHTHALIMIDE/CN

E2 1 N-(N-(3,4-DICHLOROPHENYL) CARBAMOYL) MORPHOLINE/CN

E3 0 -->

N-(N-(3-(3-METHOXY-4-HYDROXYPHENYL) PROPYL)-L-.ALPHA.-ASPARTY
L)-L-PHENYLALANINE 1-ME ESTER/CN

E4 1

N-(N-(3-AMINO-3-CARBOXYPROPYL)-3-AMINO-3-CARBOXYPROPYL) AZETI
DINE-2-CARBOXYLIC ACID/CN

E5 1 N-(N-(3-CHLOROPHENYL)-N-PHENYLAMINO) PHTHALIMIDE/CN

E6 1 N-(N-(3-PHENOXYBENZOYL) GLYCYL) GLYCINE/CN

E7 1 N-(N-(4-CARBOXYPHENYL) GLYCYL) AMINOACETONITRILE/CN

E8 1

N-(N-(5-ETHOXYCARBONYL-4-METHYL-4-PENTENOYL)-L-ALANYL)-L-VAL
INE METHYL ESTER/CN

E9 1

N-(N-(9-.BETA.-D-RIBOFURANOSYLPURIN-6-YL) CARBAMOYL) THREONINE
/CN

E10 1 N-(N-(ACRYLAMIDOMETHYL) CARBAMOYLETHYL) PHENOTHIAZINE/CN

E11 1 N-(N-(ADAMANTANE-1-CARBONYL) GLYCYL) GLYCINE/CN

E12 1 N-(N-(BENZYLOXYCARBONYL) GLYCYL) GLYCINE/CN

=>

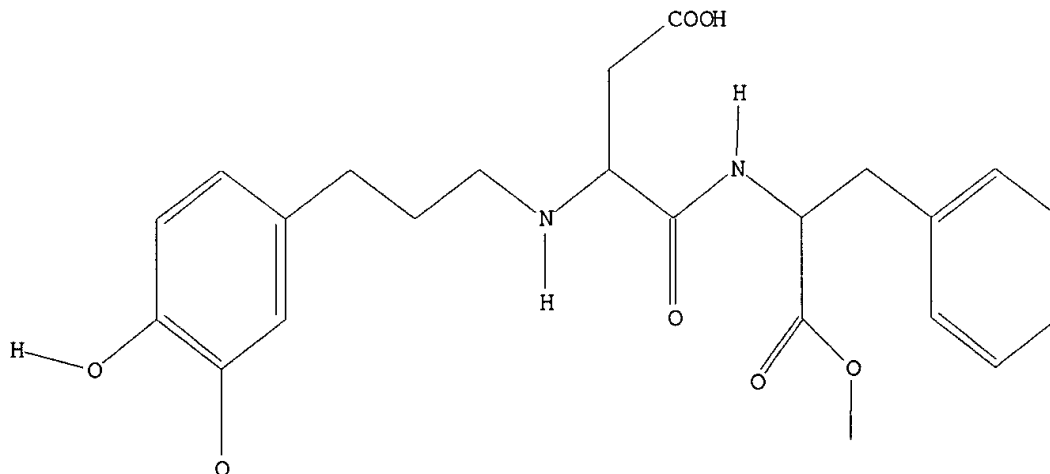
Uploading 10091500 aspartame derivative.str

L1 STRUCTURE UPLOADED

=> d 11

L1 HAS NO ANSWERS

L1 STR



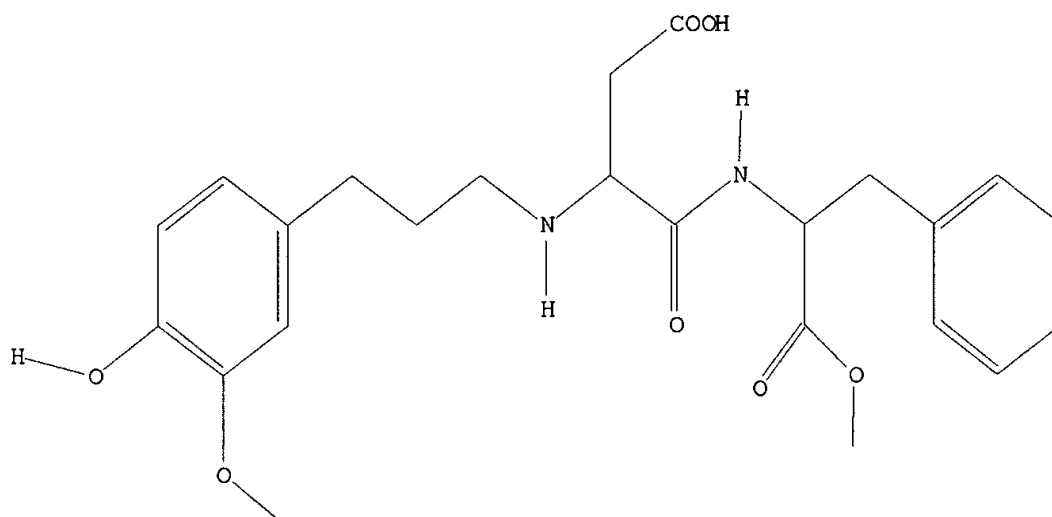
Structure attributes must be viewed using STN Express query preparation.

=>

Uploading 10091500 aspartame derivative.str

L2 STRUCTURE UPLOADED

=> d 12
L2 HAS NO ANSWERS
L2 STR



Structure attributes must be viewed using STN Express query preparation.

=> search 12 exact sam
SAMPLE SEARCH INITIATED 06:27:08 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 2 TO ITERATE

100.0% PROCESSED 2 ITERATIONS 0 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 2 TO 124
PROJECTED ANSWERS: 0 TO 0

L3 0 SEA EXA SAM L2

=> search 12 exact full
FULL SEARCH INITIATED 06:27:15 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 13 TO ITERATE

100.0% PROCESSED 13 ITERATIONS 1 ANSWERS
SEARCH TIME: 00.00.01

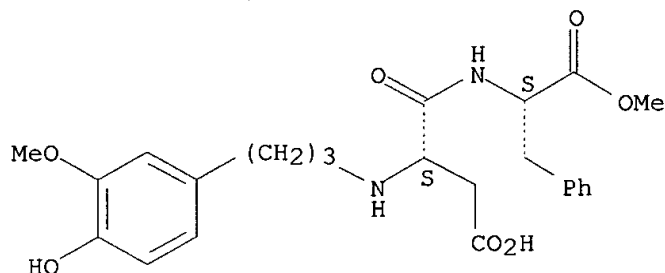
L4 1 SEA EXA FUL L2

=> d 14

L4 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2002 ACS
RN 329326-75-2 REGISTRY
CN L-Phenylalanine, N-[3-(4-hydroxy-3-methoxyphenyl)propyl]-L-.alpha.-
aspartyl-, 2-methyl ester (9CI) (CA INDEX NAME)
FS STEREOSEARCH

MF C24 H30 N2 O7
SR CA
LC STN Files: CA, CAPLUS, CASREACT

Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

2 REFERENCES IN FILE CA (1967 TO DATE)
2 REFERENCES IN FILE CAPLUS (1967 TO DATE)

=> file caplus

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

50.86

51.49

FILE 'CAPLUS' ENTERED AT 06:27:35 ON 28 AUG 2002

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FILE COVERS 1907 - 28 Aug 2002 VOL 137 ISS 9

FILE LAST UPDATED: 26 Aug 2002 (20020826/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

CAS roles have been modified effective December 16, 2001. Please check your SDI profiles to see if they need to be revised. For information on CAS roles, enter HELP ROLES at an arrow prompt or use the CAS Roles thesaurus (/RL field) in this file.

=> 14

L5

2 L4

=> d 15 1-2 ti fbib abs

L5 ANSWER 1 OF 2 CAPLUS COPYRIGHT 2002 ACS
TI Process for producing aspartyl dipeptide ester derivatives
AN 2001:833348 CAPLUS
DN 135:358168
TI Process for producing aspartyl dipeptide ester derivatives
IN Kawahara, Shigeru; Nagashima, Kazutaka; Takemoto, Tadashi
PA Ajinomoto Co., Inc., Japan
SO PCT Int. Appl., 25 pp.
CODEN: PIXXD2
DT Patent
LA Japanese
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2001085761	A1	20011115	WO 2001-JP3479	20010423
	W:				
	AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN,				
	CO, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM,				
	HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS,				
	LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO,				
	RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ,				
	VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
	RW:				
	GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY,				
	DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR, BF,				
	BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG				

JP 2000-137028 A 20000510

OS CASREACT 135:358168; MARPAT 135:358168

AB This document discloses a process for conveniently producing on an industrial scale in high yield

N-[N-[3-(phenyl)propyl]-L-.alpha.-aspartyl]-

L-phenylalanine 1-Me ester derivs., which are expected to be sweeteners, by reductively alkylating aspartame with 3-phenyl-2-propenyl aldehyde derivs. under hydrogen in the presence of a catalyst and a base.

RE.CNT 9 THERE ARE 9 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

L5 ANSWER 2 OF 2 CAPLUS COPYRIGHT 2002 ACS
TI Process for producing and purifying aspartame derivative as sweetener
AN 2001:185780 CAPLUS
DN 134:223039
TI Process for producing and purifying aspartame derivative as sweetener
IN Amino, Yusuke; Yuzawa, Kazuko; Takemoto, Tadashi
PA Ajinomoto Co., Inc., Japan
SO PCT Int. Appl., 39 pp.
CODEN: PIXXD2
DT Patent
LA Japanese
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2001018034	A1	20010315	WO 2000-JP5665	20000823
	W:				
	AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN,				
	CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR,				
	HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT,				
	LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU,				
	SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN,				
	YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				

RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY,
DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ,
CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG
JP 1999-253498 A 19990907

OS CASREACT 134:223039

AB This document discloses the following : a method for industrially
producing

N-[N-[3-(3-methoxy-4-hydroxyphenyl)propyl]-L-.alpha.-aspartyl]-L-
phenylalanine 1-Me ester which is useful as a sweetener, in particular, a
process for producing the target compd. in a high yield by the reductive
alkylation reaction of aspartame with 3-(3-methoxy-4-
hydroxyphenyl)propionaldehyde or its deriv.; a method of effectively
purifying the target compd. contaminated with impurities invading
thereinto at various prodn. stages (involving methods other than the
above-described reductive alkylation), more particularly, a method of
sepg. the target compd. in the form of highly pure crystals; the
crystals;

sweeteners contg. the same; and utilization thereof in various products
which are to be sweetened.

RE.CNT 11 THERE ARE 11 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> logoff hold

COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION

FULL ESTIMATED COST

5.37	56.86
------	-------

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE	TOTAL
ENTRY	SESSION

CA SUBSCRIBER PRICE

-1.24	-1.24
-------	-------

SESSION WILL BE HELD FOR 60 MINUTES

STN INTERNATIONAL SESSION SUSPENDED AT 06:28:46 ON 28 AUG 2002

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:sssptal623paz

PASSWORD:

* * * * * RECONNECTED TO STN INTERNATIONAL * * * * *

SESSION RESUMED IN FILE 'CAPLUS' AT 07:09:05 ON 28 AUG 2002

FILE 'CAPLUS' ENTERED AT 07:09:05 ON 28 AUG 2002

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COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION

FULL ESTIMATED COST

5.37	56.86
------	-------

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE	TOTAL
ENTRY	SESSION

CA SUBSCRIBER PRICE

-1.24	-1.24
-------	-------

=> file reg

COST IN U.S. DOLLARS

SINCE FILE	TOTAL
------------	-------

FULL ESTIMATED COST	ENTRY 5.37	SESSION 56.86
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	-1.24	-1.24

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STRUCTURE FILE UPDATES: 26 AUG 2002 HIGHEST RN 444986-65-6
 DICTIONARY FILE UPDATES: 26 AUG 2002 HIGHEST RN 444986-65-6

TSCA INFORMATION NOW CURRENT THROUGH MAY 20, 2002

Please note that search-term pricing does apply when
 conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Calculated physical property data is now available. See HELP PROPERTIES
 for more information. See STN Note 27, Searching Properties in the CAS
 Registry File, for complete details:
<http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf>

=> e 3-(4-hydroxy-3-methoxyphenyl) propionaldehyde/cn

E1 1
 3-(4-HYDROXY-3-METHOXYBENZYL)-5-HYDROXY-6,7-DIMETHOXYCHROMAN
 -4-ONE/CN

E2 1
 3-(4-HYDROXY-3-METHOXYBENZYL)-5-HYDROXY-7-METHOXYCHROMAN-4-O
 NE/CN

E3 0 --> 3-(4-HYDROXY-3-METHOXYPHENYL) PROPIONALDEHYDE/CN

E4 1 3-(4-HYDROXY-3-METHOXYPHENYL)-1,2-PROPANEDIOL/CN

E5 1 3-(4-HYDROXY-3-METHOXYPHENYL)-1,2-PROPANEDIOL TRIACETATE/CN

E6 1 3-(4-HYDROXY-3-METHOXYPHENYL)-1-PHENYL-1-PROPENE/CN

E7 1 3-(4-HYDROXY-3-METHOXYPHENYL)-1-PROPANOL/CN

E8 1 3-(4-HYDROXY-3-METHOXYPHENYL)-1-PROPENE/CN

E9 1
 3-(4-HYDROXY-3-METHOXYPHENYL)-2-(2-METHOXY-4-METHYLPHENOXY)-
 2-PROPEN-1-OL/CN

E10 1
 3-(4-HYDROXY-3-METHOXYPHENYL)-2-(2-METHOXY-4-PROPYLPHENOXY)-
 1-PROPANOL/CN

E11 1
 3-(4-HYDROXY-3-METHOXYPHENYL)-2-(2-METHOXY-5-METHYLPHENOXY)-
 2-PROPEN-1-OL/CN

E12 1
 3-(4-HYDROXY-3-METHOXYPHENYL)-2-(3,4-DIMETHOXYPHENYL)-2-PROP
 ENESULFONIC ACID/CN

=> e 3-(4-hydroxy-3-methoxyphenyl)-propionaldehyde/cn

E1 1 3-(4-HYDROXY-3-METHOXYPHENYL)-2-PROPENOIC ACID/CN

E2 1 3-(4-HYDROXY-3-METHOXYPHENYL)-5-PHENYL-1,2-DITHIOLIUM
 PERCHL
 ORATE/CN

E3 0 --> 3-(4-HYDROXY-3-METHOXYPHENYL)-PROPIONALDEHYDE/CN

E4	1	3-(4-HYDROXY-3-METHOXYPHENYL)ACROLEIN/CN
E5	1	3-(4-HYDROXY-3-METHOXYPHENYL)ACRYLIC ACID/CN
E6	1	3-(4-HYDROXY-3-METHOXYPHENYL)ALANINE/CN
E7	1	3-(4-HYDROXY-3-METHOXYPHENYL)ALLYL ALCOHOL/CN
E8	1	3-(4-HYDROXY-3-METHOXYPHENYL)LACTIC ACID/CN
E9	1	3-(4-HYDROXY-3-METHOXYPHENYL)PROPA-1,2-DIENE/CN
E10	1	3-(4-HYDROXY-3-METHOXYPHENYL)PROPANOL/CN
E11	1	3-(4-HYDROXY-3-METHOXYPHENYL)PROPIONIC ACID/CN
E12	1	3-(4-HYDROXY-3-METHOXYSTYRYL)-2,5-DIMETHYLISOXAZOLIUM
IODIDE		
		/CN

=> logoff hold

COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	1.90	58.76

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	0.00	-1.24

SESSION WILL BE HELD FOR 60 MINUTES
STN INTERNATIONAL SESSION SUSPENDED AT 07:11:53 ON 28 AUG 2002

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:sssptal623paz

PASSWORD:

* * * * * RECONNECTED TO STN INTERNATIONAL * * * * *
SESSION RESUMED IN FILE 'REGISTRY' AT 07:17:02 ON 28 AUG 2002
FILE 'REGISTRY' ENTERED AT 07:17:02 ON 28 AUG 2002
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COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	1.90	58.76
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	0.00	-1.24

=>

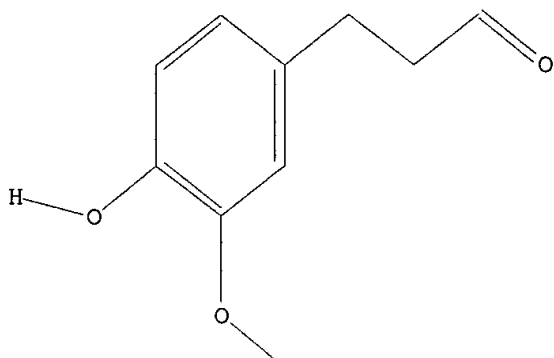
Uploading 10091500 aldehyde.str

L6 STRUCTURE UPLOADED

=> d l6

L6 HAS NO ANSWERS

L6 STR



Structure attributes must be viewed using STN Express query preparation.

=> search l6 exact full

FULL SEARCH INITIATED 07:18:00 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 124 TO ITERATE

100.0% PROCESSED 124 ITERATIONS

1 ANSWERS

SEARCH TIME: 00.00.01

L7 1 SEA EXA FUL L6

=> d 17

L7 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2002 ACS

RN 80638-48-8 REGISTRY

CN Benzenepropanal, 4-hydroxy-3-methoxy- (9CI) (CA INDEX NAME)

OTHER NAMES:

CN 3-(4'-Hydroxy-3'-methoxyphenyl)-1-propanal

CN 4-Propanal-2-methoxyphenol

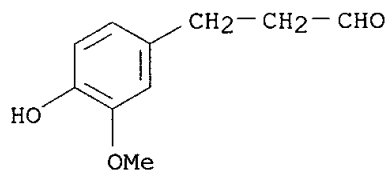
FS 3D CONCORD

MF C10 H12 O3

LC STN Files: CA, CAPLUS, CASREACT, CHEMLIST, TOXCENTER

Other Sources: DSL**, EINECS**

(**Enter CHEMLIST File for up-to-date regulatory information)



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

11 REFERENCES IN FILE CA (1967 TO DATE)

11 REFERENCES IN FILE CAPLUS (1967 TO DATE)

=> file caplus

COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	51.24	108.10
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	0.00	-1.24

FILE 'CAPLUS' ENTERED AT 07:18:17 ON 28 AUG 2002
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FILE COVERS 1907 - 28 Aug 2002 VOL 137 ISS 9
FILE LAST UPDATED: 26 Aug 2002 (20020826/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

CAS roles have been modified effective December 16, 2001. Please check your SDI profiles to see if they need to be revised. For information on CAS roles, enter HELP ROLES at an arrow prompt or use the CAS Roles thesaurus (/RL field) in this file.

=> 17

L8 11 L7

=> save temp l8 sweetaldehy/a
ANSWER SET L8 HAS BEEN SAVED AS 'SWEETALDEH/A'

=>

COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	0.40	108.50
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	0.00	-1.24

FILE 'STNGUIDE' ENTERED AT 07:19:02 ON 28 AUG 2002
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AND TECHNOLOGY CORPORATION, AND FACHINFORMATIONSZENTRUM KARLSRUHE

FILE CONTAINS CURRENT INFORMATION.
LAST RELOADED: Aug 23, 2002 (20020823/UP).

=>

NAME	CREATED	NOTES/TITLE
ADIACRYST/A	TEMP	1001 ANSWERS IN FILE CAPLUS
ALKYLATIN/L	13 DEC 2001	9 L-NUMBERS
ASPARTAME/A	TEMP	2346 ANSWERS IN FILE CAPLUS
A408SRCH/L	TEMP	42 L-NUMBERS
CONIFERAL/A	TEMP	388 ANSWERS IN FILE CAPLUS
GLYPHSRCH/L	TEMP	13 L-NUMBERS
INDIUMCL3/A	30 MAY 2001	1 ANSWER IN FILE REGISTRY
LTWENTAUGFOR/A	04 AUG 2001	72 ANSWERS IN FILE CAPLUS
NASTATINS/A	TEMP	144 ANSWERS IN FILE CAPLUS
NEOTAMECRYST/A	24 APR 2001	59 ANSWERS IN FILE CAPLUS
NVLARMFULGEN/A	19 APR 2001	196 ANSWERS IN FILE REGISTRY
POHBENZALDEH/A	10 JUL 2001	5519 ANSWERS IN FILE CAPLUS
PROCTYLCMPD/A	TEMP	10 ANSWERS IN FILE CAPLUS
PROCTYLSRCH/L	TEMP	4 L-NUMBERS
PROSTACMPD15/A	01 AUG 2001	34 ANSWERS IN FILE CAPLUS
STILLEAPP/L	07 JAN 2002	17 L-NUMBERS
SWEET/L	TEMP	6 L-NUMBERS
SWEETALDEH/A	TEMP	11 ANSWERS IN FILE CAPLUS
TWOAMINOPOLY/Q	16 APR 2001	UPLOADED STRUCTURE

=>

NO SAVED SDI REQUESTS

=>

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.06	108.56
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	0.00	-1.24

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FILE COVERS 1907 - 28 Aug 2002 VOL 137 ISS 9
FILE LAST UPDATED: 26 Aug 2002 (20020826/ED)

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=>

L9 (1)SEA FILE=REGISTRY ABB=ON PLU=ON ASPARTAME/CN
L10 2346 SEA FILE=CAPLUS ABB=ON PLU=ON L9

=>

L11 (1)SEA FILE=REGISTRY ABB=ON PLU=ON CONIFERALDEHYDE/CN
L12 388 SEA FILE=CAPLUS ABB=ON PLU=ON L11

=> file caplus

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.40	108.96

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	0.00	-1.24

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=> d his

(FILE 'HOME' ENTERED AT 06:22:05 ON 28 AUG 2002)

FILE 'REGISTRY' ENTERED AT 06:23:52 ON 28 AUG 2002

E
N-[N-[3-(3-METHOXY-4-HYDROXYPHENYL)PROPYL]-L-.ALPHA.-ASPARTYL
E
N-(N-(3-(3-METHOXY-4-HYDROXYPHENYL)PROPYL)-L-.ALPHA.-ASPARTYL
L1 STRUCTURE UPLOADED

L2 STRUCTURE UPLOADED
L3 0 SEARCH L2 EXACT SAM
L4 1 SEARCH L2 EXACT FULL

FILE 'CAPLUS' ENTERED AT 06:27:35 ON 28 AUG 2002
L5 2 L4

FILE 'REGISTRY' ENTERED AT 07:09:10 ON 28 AUG 2002
E 3-(4-HYDROXY-3-METHOXYPHENYL) PROPIONALDEHYDE/CN
E 3-(4-HYDROXY-3-METHOXYPHENYL)-PROPIONALDEHYDE/CN
L6 STRUCTURE UPLOADED
L7 1 SEARCH L6 EXACT FULL

FILE 'CAPLUS' ENTERED AT 07:18:17 ON 28 AUG 2002
L8 11 L7
SAVE TEMP L8 SWEETALDEH/A

FILE 'STNGUIDE' ENTERED AT 07:19:02 ON 28 AUG 2002

FILE 'CAPLUS' ENTERED AT 07:19:47 ON 28 AUG 2002
ACT ASPARTAME/A

L9 (1)SEA FILE=REGISTRY ABB=ON PLU=ON ASPARTAME/CN
L10 2346 SEA FILE=CAPLUS ABB=ON PLU=ON L9

ACT CONIFERAL/A

L11 (1)SEA FILE=REGISTRY ABB=ON PLU=ON CONIFERALDEHYDE/CN
L12 388 SEA FILE=CAPLUS ABB=ON PLU=ON L11

FILE 'CAPLUS' ENTERED AT 07:20:01 ON 28 AUG 2002

=> l8 and l10
L13 1 L8 AND L10

=> d l13 ti fbib abs

L13 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2002 ACS
TI Process for producing and purifying aspartame derivative as sweetener
AN 2001:185780 CAPLUS
DN 134:223039
TI Process for producing and purifying aspartame derivative as sweetener
IN Amino, Yusuke; Yuzawa, Kazuko; Takemoto, Tadashi
PA Ajinomoto Co., Inc., Japan
SO PCT Int. Appl., 39 pp.
CODEN: PIXXD2
DT Patent
LA Japanese
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2001018034	A1	20010315	WO 2000-JP5665	20000823
W:	AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				

RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY,
DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ,
CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG

JP 1999-253498 A 19990907

OS CASREACT 134:223039

AB This document discloses the following : a method for industrially
producing

N-[N-[3-(3-methoxy-4-hydroxyphenyl)propyl]-L-.alpha.-aspartyl]-L-
phenylalanine 1-Me ester which is useful as a sweetener, in particular, a
process for producing the target compd. in a high yield by the reductive
alkylation reaction of aspartame with 3-(3-methoxy-4-
hydroxyphenyl)propionaldehyde or its deriv.; a method of effectively
purifying the target compd. contaminated with impurities invading
thereinto at various prodn. stages (involving methods other than the
above-described reductive alkylation), more particularly, a method of
sepg. the target compd. in the form of highly pure crystals; the
crystals;

sweeteners contg. the same; and utilization thereof in various products
which are to be sweetened.

RE.CNT 11 THERE ARE 11 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> l8 and l12

L14 6 L8 AND L12

=> d l14 1-6 ti

L14 ANSWER 1 OF 6 CAPLUS COPYRIGHT 2002 ACS

TI Process for producing and purifying aspartame derivative as sweetener

L14 ANSWER 2 OF 6 CAPLUS COPYRIGHT 2002 ACS

TI Retention of lignin in seagrasses: angiosperms that returned to the sea

L14 ANSWER 3 OF 6 CAPLUS COPYRIGHT 2002 ACS

TI Chemical composition of alkaline pulping products of lignin model
compounds

L14 ANSWER 4 OF 6 CAPLUS COPYRIGHT 2002 ACS

TI Thermolytic decomposition of coniferyl alcohol

L14 ANSWER 5 OF 6 CAPLUS COPYRIGHT 2002 ACS

TI Ultrafiltration and pyrolysis gas chromatography mass spectrometry of
chlorolignins in pulp mill effluent

L14 ANSWER 6 OF 6 CAPLUS COPYRIGHT 2002 ACS

TI Topical cosmetics containing 1,7-diphenyl-4-hepten-3-one for skin
disorder
treatment

=> logoff hold

COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
13.10	122.06

FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE	TOTAL
ENTRY	SESSION
-0.62	-1.86

CA SUBSCRIBER PRICE

SESSION WILL BE HELD FOR 60 MINUTES
STN INTERNATIONAL SESSION SUSPENDED AT 07:33:50 ON 28 AUG 2002

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:sssptal623paz

PASSWORD:

* * * * * RECONNECTED TO STN INTERNATIONAL * * * * *
SESSION RESUMED IN FILE 'CAPLUS' AT 07:51:23 ON 28 AUG 2002
FILE 'CAPLUS' ENTERED AT 07:51:23 ON 28 AUG 2002
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COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	13.10	122.06

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	-0.62	-1.86

=> file reg

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	13.89	122.85

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	-0.62	-1.86

FILE 'REGISTRY' ENTERED AT 07:52:26 ON 28 AUG 2002
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STRUCTURE FILE UPDATES: 26 AUG 2002 HIGHEST RN 444986-65-6
DICTIONARY FILE UPDATES: 26 AUG 2002 HIGHEST RN 444986-65-6

TSCA INFORMATION NOW CURRENT THROUGH MAY 20, 2002

Please note that search-term pricing does apply when
conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Calculated physical property data is now available. See HELP PROPERTIES
for more information. See STN Note 27, Searching Properties in the CAS
Registry File, for complete details:
<http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf>

=> e Benzenepropanal, 4-hydroxy-3-methoxy-,cn

E1 106272 BENZENEPROPAN/BI
E2 2540 BENZENEPROPANAL/BI
E3 0 --> BENZENEPROPANAL, 4-HYDROXY-3-METHOXY-,CN/BI

E4	60	BENZENEPROPANALATO/BI
E5	1	BENZENEPROPANALDEHYDE/BI
E6	538	BENZENEPROPANAMID/BI
E7	536	BENZENEPROPANAMIDATO/BI
E8	24838	BENZENEPROPANAMIDE/BI
E9	1	BENZENEPROPANAMIDO/BI
E10	574	BENZENEPROPANAMIN/BI
E11	22	BENZENEPROPANAMINATO/BI
E12	5004	BENZENEPROPANAMINE/BI

=> e 4-hydroxy-3-methoxy-Benzenepropanal cn

E1	1	4-AMINO-2,6-PYRIMIDINODITHIOCARBAMATE/BI
E2	1	4-HYDROXY-2-METHYL-N-2-PYRIDINYL-2H-1,2-BENZOTHAZINE-3-CARB OXAMIDE 1,1-DIOXIDE-NN/BI
E3	0 -->	4-HYDROXY-3-METHOXY-BENZENEPROPANAL CN/BI
E4	3	4-ISOCYANATOBENZENE/BI
E5	1	4-METHOXY-.ALPHA.-METHYLBENZENEACETATE/BI
E6	2	4-METHYL-2-OXO-2H-1-BENZOPYRAN-7-YL/BI
E7	1	4-METHYLBENZENESULFONAMIDATO/BI
E8	1	4-METHYLBENZENESULFONATO-.KAPPA.O/BI
E9	1	4-METHYLPHENYL 6-DIAZO-5,6-DIHYDRO-5-OXO-1-NAPHTHALENESULFON ATE/BI
E10	1	4-MORPHOLINO-1,3,5-TRIAZINE-2,6-DIYL/BI
E11	2	4-NITROPHENYL/BI
E12	1	4-O-.BETA.-D-GALACTOPYRANOSYL-D-GLUCONATO-O1,O2,O3/BI

=> file caplus

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	1.14	123.99
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	0.00	-1.86

FILE 'CAPLUS' ENTERED AT 07:54:07 ON 28 AUG 2002
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=> d his

(FILE 'HOME' ENTERED AT 06:22:05 ON 28 AUG 2002)

FILE 'REGISTRY' ENTERED AT 06:23:52 ON 28 AUG 2002

E
N-[N-[3-(3-METHOXY-4-HYDROXYPHENYL) PROPYL]-L-.ALPHA.-ASPARTYL
E
N-(N-(3-(3-METHOXY-4-HYDROXYPHENYL) PROPYL)-L-.ALPHA.-ASPARTYL
L1 STRUCTURE UPLOADED
L2 STRUCTURE UPLOADED
L3 0 SEARCH L2 EXACT SAM
L4 1 SEARCH L2 EXACT FULL

FILE 'CAPLUS' ENTERED AT 06:27:35 ON 28 AUG 2002

L5 2 L4

FILE 'REGISTRY' ENTERED AT 07:09:10 ON 28 AUG 2002

E 3-(4-HYDROXY-3-METHOXYPHENYL) PROPIONALDEHYDE/CN
E 3-(4-HYDROXY-3-METHOXYPHENYL)-PROPIONALDEHYDE/CN
L6 STRUCTURE UPLOADED
L7 1 SEARCH L6 EXACT FULL

FILE 'CAPLUS' ENTERED AT 07:18:17 ON 28 AUG 2002

L8 11 L7
SAVE TEMP L8 SWEETALDEH/A

FILE 'STNGUIDE' ENTERED AT 07:19:02 ON 28 AUG 2002

FILE 'CAPLUS' ENTERED AT 07:19:47 ON 28 AUG 2002
ACT ASPARTAME/A

L9 (1)SEA FILE=REGISTRY ABB=ON PLU=ON ASPARTAME/CN
L10 2346 SEA FILE=CAPLUS ABB=ON PLU=ON L9

ACT CONIFERAL/A

L11 (1)SEA FILE=REGISTRY ABB=ON PLU=ON CONIFERALDEHYDE/CN
L12 388 SEA FILE=CAPLUS ABB=ON PLU=ON L11

FILE 'CAPLUS' ENTERED AT 07:20:01 ON 28 AUG 2002

L13 1 L8 AND L10
L14 6 L8 AND L12

FILE 'REGISTRY' ENTERED AT 07:52:26 ON 28 AUG 2002

E BENZENEPROPANAL, 4-HYDROXY-3-METHOXY-,CN
E 4-HYDROXY-3-METHOXY-BENZENEPROPANAL CN

FILE 'CAPLUS' ENTERED AT 07:54:07 ON 28 AUG 2002

=> reg

915 REG
41 REGS

L15 952 REG
(REG OR REGS)

=> file reg

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	2.02	126.01

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	0.00	-1.86

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DICTIONARY FILE UPDATES: 26 AUG 2002 HIGHEST RN 444986-65-6

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for more information. See STNnote 27, Searching Properties in the CAS
Registry File, for complete details:
<http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf>

=> e coniferaldehyde/cn

E1 1 CONIDIOSPORE SURFACE PROTEIN (TRICHODERMA HARZIANUM STRAIN
A TCC 32173 GENE CMP1)/CN
E2 1 CONIFEGEROL/CN
E3 1 --> CONIFERALDEHYDE/CN
E4 1 CONIFERALDEHYDE METHOXYMETHYL ETHER/CN
E5 1 CONIFERALDEHYDE METHYL ETHER/CN
E6 1 CONIFERAN/CN
E7 1 CONIFERIC ACID/CN
E8 1 CONIFERIN/CN
E9 1 CONIFERIN .BETA.-GLUCOSIDASE/CN
E10 1 CONIFERIN PENTAACETATE/CN
E11 1 CONIFERIN, DIHYDRATE/CN
E12 1 CONIFERIN-HYDROLYZING .BETA.-GLUCOSIDASE/CN

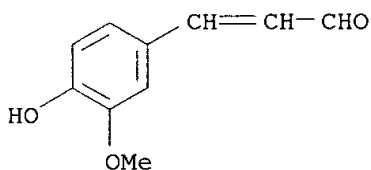
=> e3

L16 1 CONIFERALDEHYDE/CN

=> d 116

L16 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2002 ACS
RN 458-36-6 REGISTRY
CN 2-Propenal, 3-(4-hydroxy-3-methoxyphenyl)- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Cinnamaldehyde, 4-hydroxy-3-methoxy- (6CI, 8CI)
OTHER NAMES:

CN 2-Methoxy-4-(3-oxo-1-propenyl)phenol
 CN 3-(4-Hydroxy-3-methoxyphenyl)acrolein
 CN 3-Methoxy-4-hydroxycinnamaldehyde
 CN 4-Hydroxy-3-methoxycinnamaldehyde
 CN **Coniferaldehyde**
 CN Coniferyl aldehyde
 CN Ferulaldehyde
 CN Ferulyl aldehyde
 CN p-Coniferaldehyde
 FS 3D CONCORD
 MF C10 H10 O3
 CI COM
 LC STN Files: AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS,
 BIOTECHNO, CA, CAOLD, CAPLUS, CASREACT, CHEMCATS, CHEMLIST, CSCHEM,
 DDFU, DRUGU, EMBASE, HODOC*, MEDLINE, NAPRALERT, NIOSHTIC, PIRA,
 SPECINFO, TOXCENTER, USPATFULL
 (*File contains numerically searchable property data)
 Other Sources: EINECS**
 (**Enter CHEMLIST File for up-to-date regulatory information)



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

388 REFERENCES IN FILE CA (1967 TO DATE)
 3 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
 388 REFERENCES IN FILE CAPLUS (1967 TO DATE)
 28 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

=> file capls
 'CAPLS' IS NOT A VALID FILE NAME
 SESSION CONTINUES IN FILE 'REGISTRY'
 Enter "HELP FILE NAMES" at an arrow prompt (=>) for a list of files
 that are available. If you have requested multiple files, you can
 specify a corrected file name or you can enter "IGNORE" to continue
 accessing the remaining file names entered.

=> file caplus		
COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	5.96	131.97
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	0.00	-1.86

FILE 'CAPLUS' ENTERED AT 07:55:19 ON 28 AUG 2002
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=> d his

(FILE 'HOME' ENTERED AT 06:22:05 ON 28 AUG 2002)

FILE 'REGISTRY' ENTERED AT 06:23:52 ON 28 AUG 2002

E

N-[N-(3-(3-METHOXY-4-HYDROXYPHENYL) PROPYL)-L-.ALPHA.-ASPARTYL

E

N-(N-(3-(3-METHOXY-4-HYDROXYPHENYL) PROPYL)-L-.ALPHA.-ASPARTYL

L1 STRUCTURE UPLOADED

L2 STRUCTURE UPLOADED

L3 0 SEARCH L2 EXACT SAM

L4 1 SEARCH L2 EXACT FULL

FILE 'CAPLUS' ENTERED AT 06:27:35 ON 28 AUG 2002

L5 2 L4

FILE 'REGISTRY' ENTERED AT 07:09:10 ON 28 AUG 2002

E 3-(4-HYDROXY-3-METHOXYPHENYL) PROPIONALDEHYDE/CN

E 3-(4-HYDROXY-3-METHOXYPHENYL)-PROPIONALDEHYDE/CN

L6 STRUCTURE UPLOADED

L7 1 SEARCH L6 EXACT FULL

FILE 'CAPLUS' ENTERED AT 07:18:17 ON 28 AUG 2002

L8 11 L7

SAVE TEMP L8 SWEETALDEH/A

FILE 'STNGUIDE' ENTERED AT 07:19:02 ON 28 AUG 2002

FILE 'CAPLUS' ENTERED AT 07:19:47 ON 28 AUG 2002

ACT ASPARTAME/A

L9 (1)SEA FILE=REGISTRY ABB=ON PLU=ON ASPARTAME/CN

L10 2346 SEA FILE=CAPLUS ABB=ON PLU=ON L9

ACT CONIFERAL/A

L11 (1)SEA FILE=REGISTRY ABB=ON PLU=ON CONIFERALDEHYDE/CN
L12 388 SEA FILE=CAPLUS ABB=ON PLU=ON L11

FILE 'CAPLUS' ENTERED AT 07:20:01 ON 28 AUG 2002
L13 1 L8 AND L10
L14 6 L8 AND L12

FILE 'REGISTRY' ENTERED AT 07:52:26 ON 28 AUG 2002
E BENZENEPROPANAL, 4-HYDROXY-3-METHOXY-,CN
E 4-HYDROXY-3-METHOXY-BENZENEPROPANAL CN

FILE 'CAPLUS' ENTERED AT 07:54:07 ON 28 AUG 2002
L15 952 REG

FILE 'REGISTRY' ENTERED AT 07:54:37 ON 28 AUG 2002
E CONIFERALDEHYDE/CN
L16 1 E3

FILE 'CAPLUS' ENTERED AT 07:55:19 ON 28 AUG 2002

=> rhodium
52919 RHODIUM
30 RHODIUMS
L17 52920 RHODIUM
(RHODIUM OR RHODIUMS)

=> l12 and l17
L18 1 L12 AND L17

=> d l18 ti

L18 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2002 ACS
TI Process for producing and purifying aspartame derivative as sweetener

=> logoff hold		
COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	2.70	134.67
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	0.00	-1.86

SESSION WILL BE HELD FOR 60 MINUTES
STN INTERNATIONAL SESSION SUSPENDED AT 07:56:45 ON 28 AUG 2002

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:sssptal623paz